Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080515\
Data File : BE090378.D
Acq On : 6 Aug 2015 6:08
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 06 13:38:45 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 02:22:05 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	108	0.00
2		1,4-Dioxane	1.000	1.111	-11.1	106	0.00
3		n-Nitrosodimethylamine	1.000	1.001	-0.1	109	0.00
4	S	2-Fluorophenol	1.000	1.055	-5.5	119	0.00
5	S	Phenol-d6	1.000	0.960	4.0	112	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	108	0.00
7	S	Nitrobenzene-d5	1.000	0.947	5.3	109	0.00
8		Nitrobenzene	1.000	0.904	9.6	109	0.00
9		Naphthalene	1.000	1.002	-0.2	109	0.00
10		Hexachlorobutadiene	1.000	1.107	-10.7	109	0.00
11		2-Methylnaphthalene	1.000	1.020	-2.0	115	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	108	0.00
13	S	2,4,6-Tribromophenol	1.000	1.023	-2.3	122	0.00
14	S	2-Fluorobiphenyl	1.000	1.053	-5.3	112	0.00
15		Acenaphthylene	1.000	1.025	-2.5	111	0.00
16		Acenaphthene	1.000	1.007	-0.7	108	0.00
17		Fluorene	1.000	1.025	-2.5	109	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	104	0.00
19		4-Bromophenyl-phenylether	1.000	1.049	-4.9	111	0.00
20		Hexachlorobenzene	1.000	1.023	-2.3	106	0.00
21		Pentachlorophenol	1.000	1.277	-27.7#	169	0.00
22		Phenanthrene	1.000	1.021	-2.1	107	0.00
23		Anthracene	1.000	1.014	-1.4	108	-0.02
24		Fluoranthene	1.000	1.035	-3.5	111	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	108	0.00
26		Pyrene	1.000	1.017	-1.7	114	0.00
27	S	Terphenyl-d14	1.000	1.025	-2.5	112	0.00
28		Benzo(a)anthracene	1.000	0.992	0.8	119	0.00
29		Chrysene	1.000	0.971	2.9	102	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.901	9.9	108	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.921	7.9	118	0.00
32	I	Perylene-d12	5.000	5.000	0.0	109	0.00
33		Benzo(b)fluoranthene	1.000	1.027	-2.7	123	0.00
34		Benzo(k)fluoranthene	1.000	0.978	2.2	102	0.00
35	C	Benzo(a)pyrene	1.000	0.986	1.4	116	0.00
36		Dibenzo(a,h)anthracene	1.000	0.951	4.9	121	0.00
37		Benzo(g,h,i)perylene	1.000	1.007	-0.7	113	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0